

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092618\
 Data File : BF109334.D
 Acq On : 26 Sep 2018 14:17
 Operator : JU/SJ
 Sample : J5064-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.228	359	364	373	rBV	530498	631863	25.84%	2.745%
2	4.881	470	475	487	rVB	557722	614550	25.13%	2.670%
3	5.310	544	548	551	rBV	1406055	1238026	50.63%	5.379%
4	6.339	719	723	726	rBV	1682361	1571050	64.25%	6.826%
5	6.475	742	746	749	rBV	1771513	1471437	60.18%	6.393%
6	6.704	782	785	789	rBV	763834	618632	25.30%	2.688%
7	6.781	794	798	801	rBV	832310	681220	27.86%	2.960%
8	6.863	808	812	815	rBV	1782983	1406276	57.51%	6.110%
9	7.263	876	880	883	rBV	1162988	1096881	44.86%	4.766%
10	7.986	1000	1003	1009	rBV	986368	773308	31.62%	3.360%
11	8.875	1151	1154	1156	rBV	128030	123698	5.06%	0.537%
12	8.957	1165	1168	1171	rBV	167457	151738	6.21%	0.659%
13	8.998	1172	1175	1178	rVV	163272	173653	7.10%	0.755%
14	9.027	1178	1180	1182	rVB	167763	125022	5.11%	0.543%
15	9.063	1182	1186	1189	rBV	2303325	1929850	78.92%	8.385%
16	9.157	1198	1202	1206	rBV2	385365	545706	22.32%	2.371%
17	9.216	1208	1212	1214	rBV4	231553	334794	13.69%	1.455%
18	9.239	1214	1216	1218	rVB	140948	86981	3.56%	0.378%
19	9.274	1220	1222	1225	rVB	146867	111936	4.58%	0.486%
20	9.310	1225	1228	1232	rBV3	193459	303266	12.40%	1.318%
21	9.351	1232	1235	1237	rBV	186680	188592	7.71%	0.819%
22	9.380	1237	1240	1242	rBV2	202342	268987	11.00%	1.169%
23	9.416	1243	1246	1249	rBV3	446825	590528	24.15%	2.566%
24	9.457	1251	1253	1254	rBV	399396	277435	11.35%	1.205%
25	9.486	1255	1258	1260	rVB	513103	433880	17.74%	1.885%
26	9.545	1266	1268	1270	rVB2	368523	272503	11.14%	1.184%
27	9.569	1270	1272	1274	rBV2	237626	196960	8.05%	0.856%
28	9.627	1277	1282	1283	rBV4	471796	678497	27.75%	2.948%
29	9.663	1285	1288	1291	rVB2	715779	681349	27.86%	2.960%
30	9.698	1291	1294	1296	rBV2	722376	749980	30.67%	3.259%
31	9.739	1297	1301	1307	rBV3	1361438	2445253	100.00%	10.625%
32	10.027	1347	1350	1353	rBV	1053524	1185395	48.48%	5.151%
33	15.256	2234	2239	2242	rVV	603957	715442	29.26%	3.109%
34	15.292	2242	2245	2251	rVB	70505	96920	3.96%	0.421%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	16.151	2388	2391	2399	rvB3	46937	84416	3.45%	0.367%
36	17.509	2616	2622	2630	rvB2	64808	158859	6.50%	0.690%

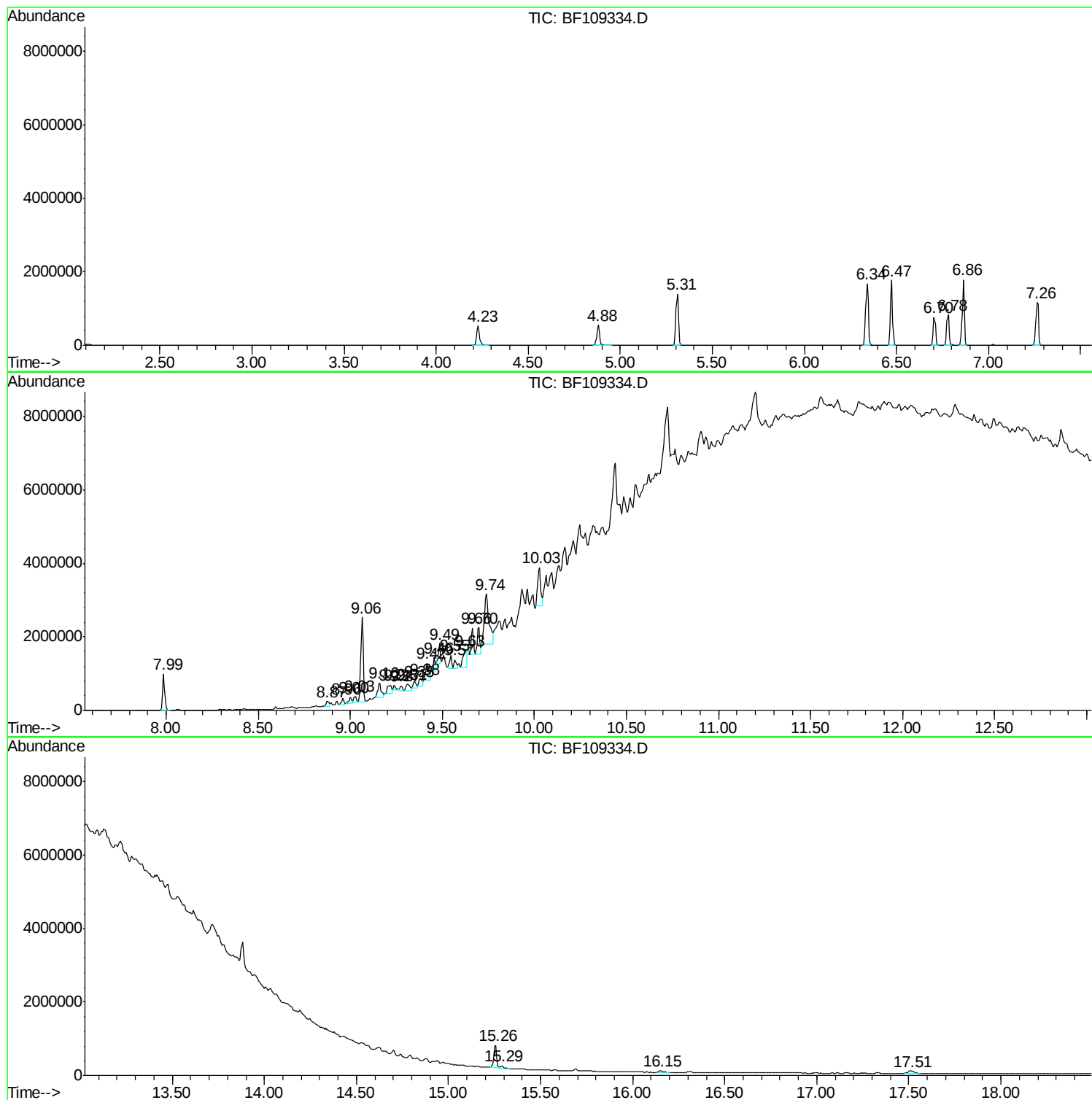
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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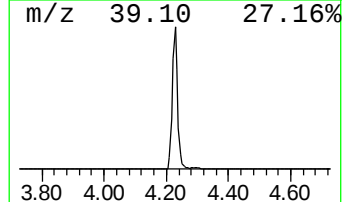
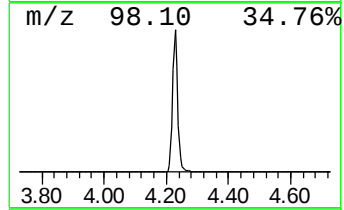
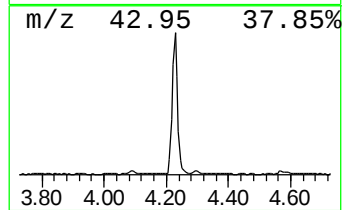
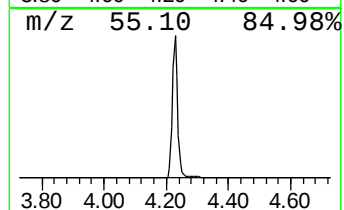
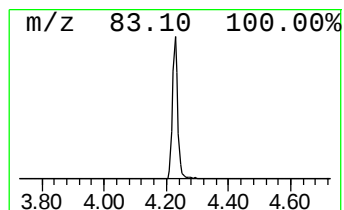
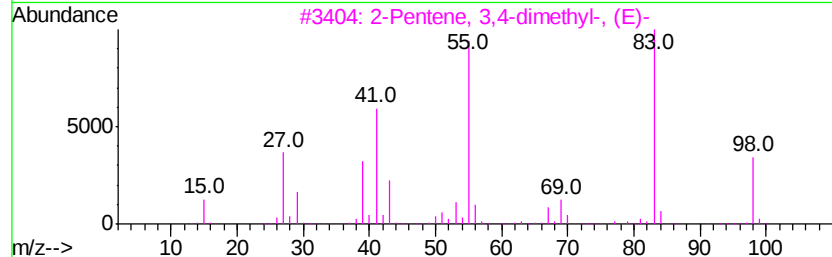
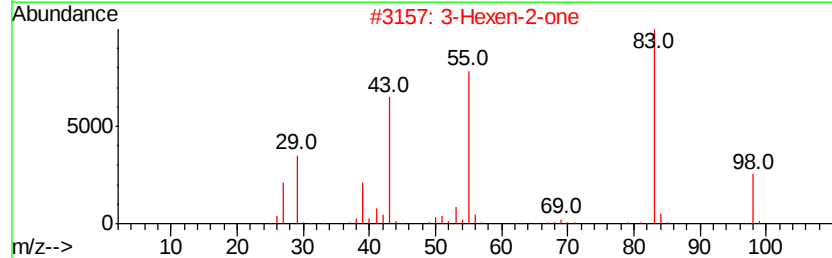
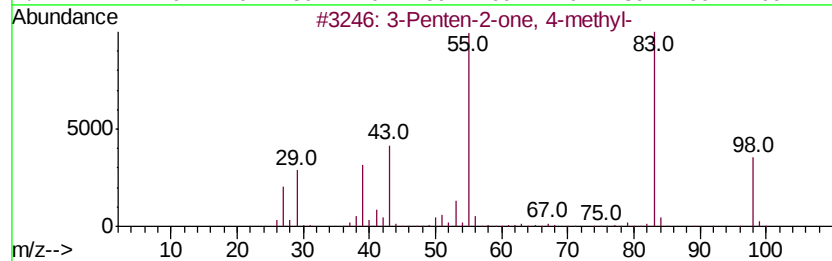
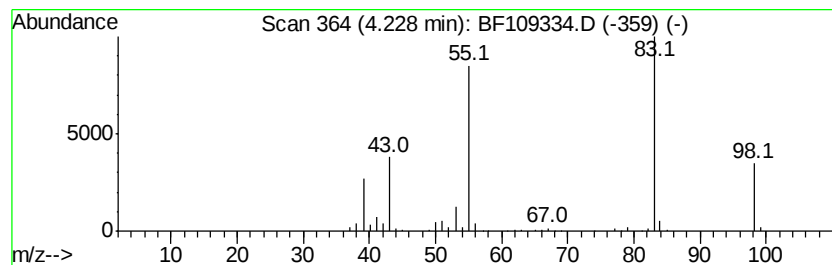
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.23	20.43 ng	631863	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80



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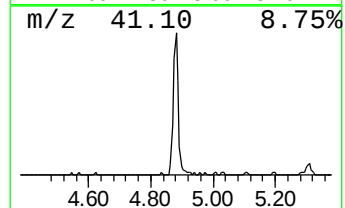
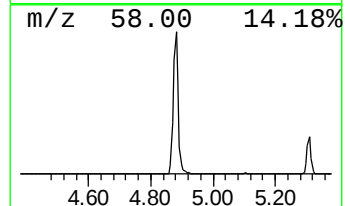
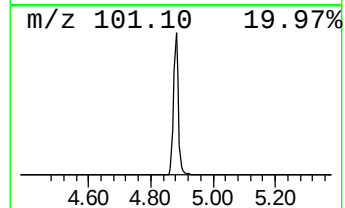
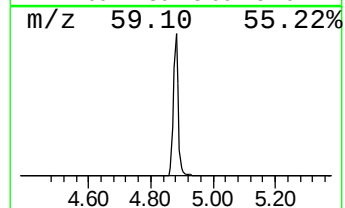
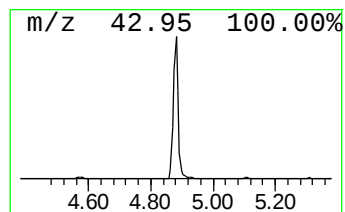
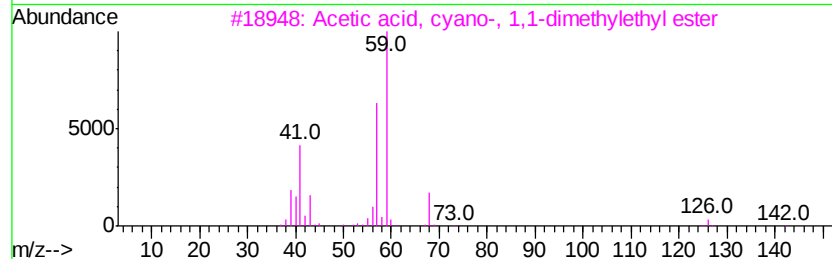
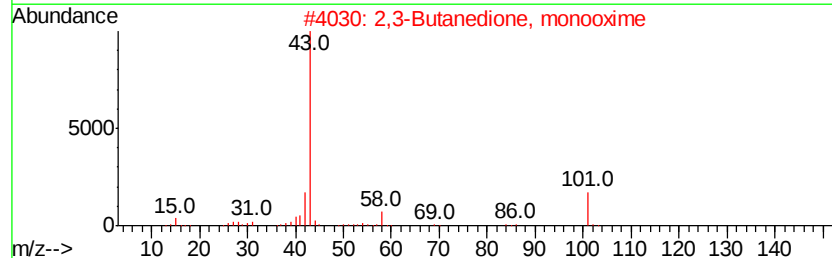
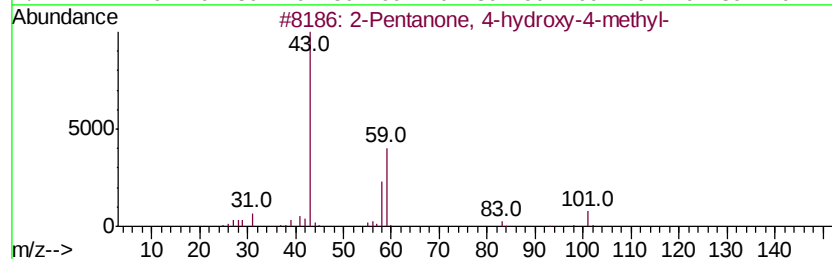
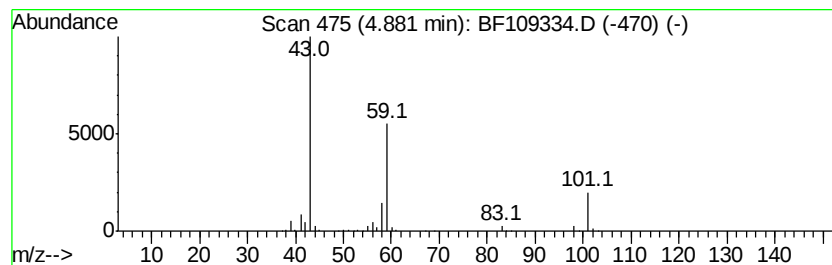
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	19.87 ng	614550	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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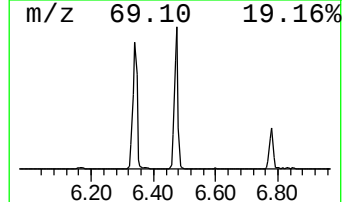
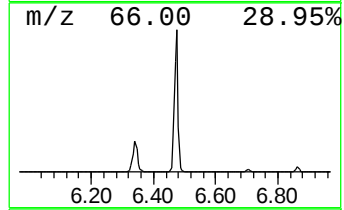
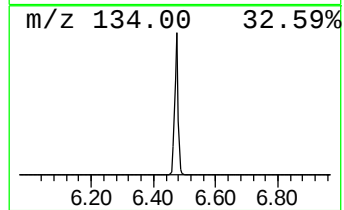
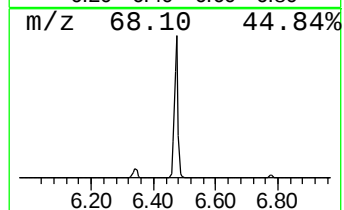
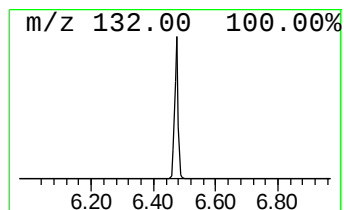
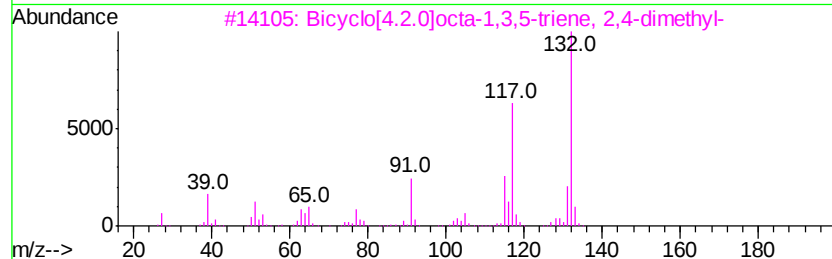
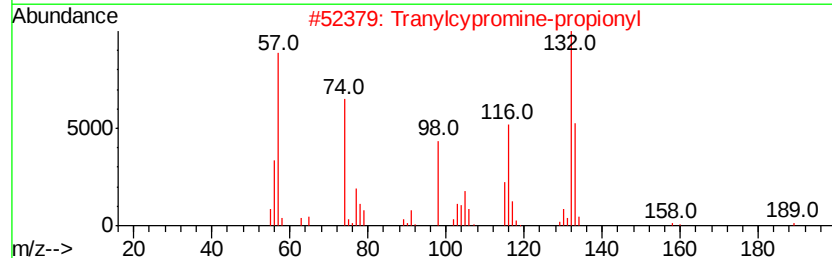
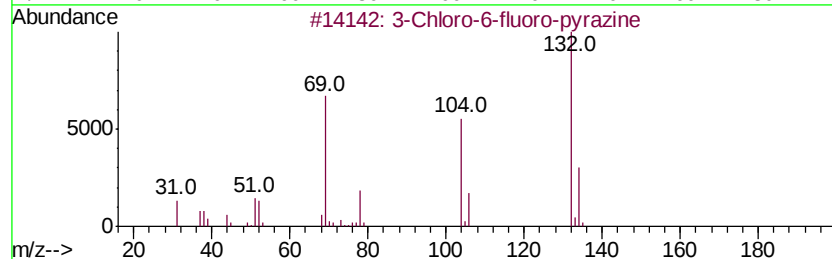
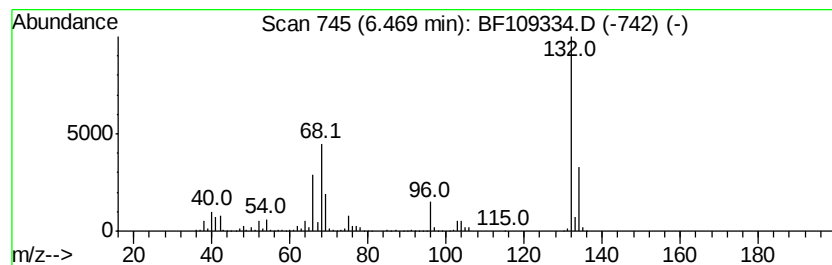
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	47.57 ng	1471440	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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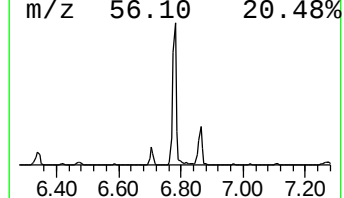
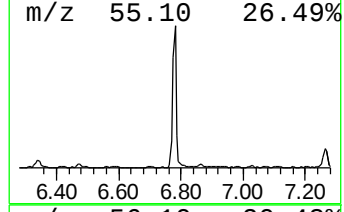
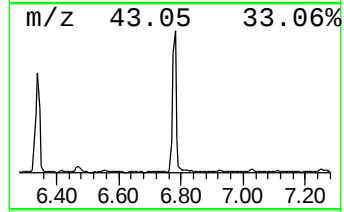
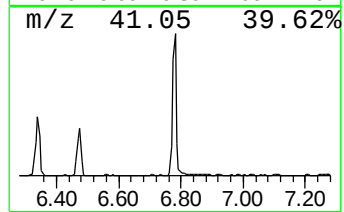
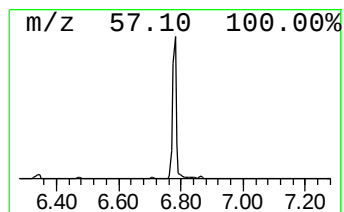
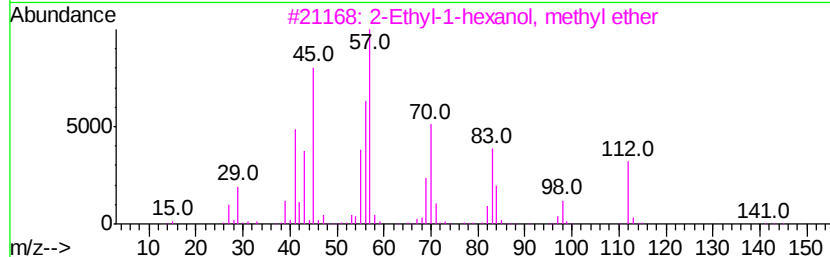
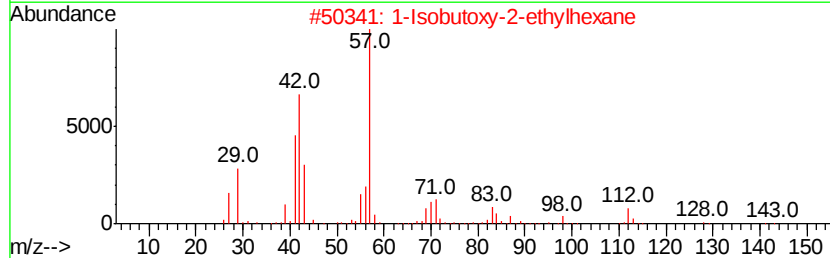
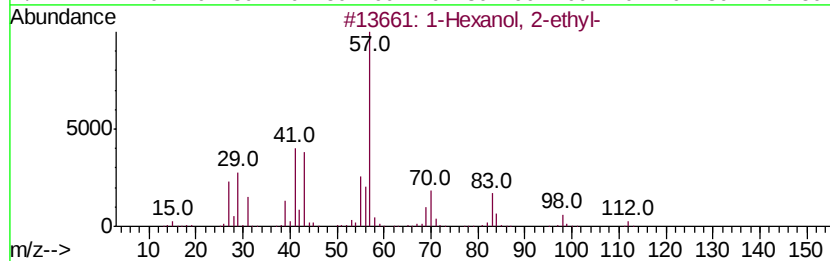
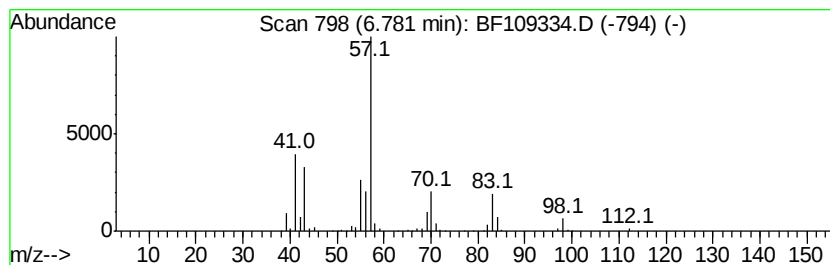
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.78	22.02 ng	681220	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2		1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	50
3		2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	50
4		Oxalic acid, isobutyl octyl ester	258	C14H26O4	1000309-37-3	47
5		Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	45



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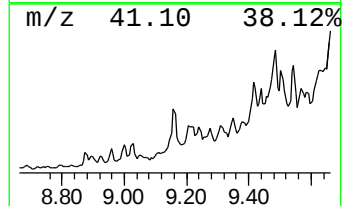
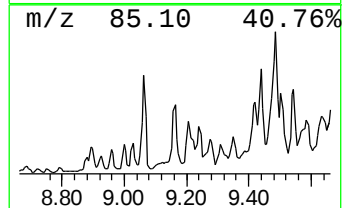
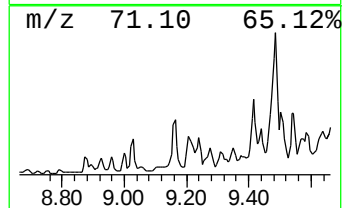
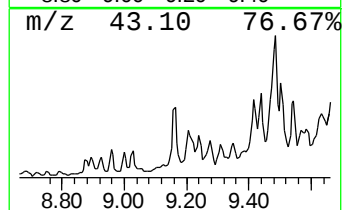
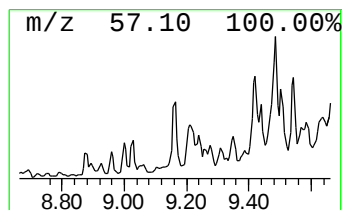
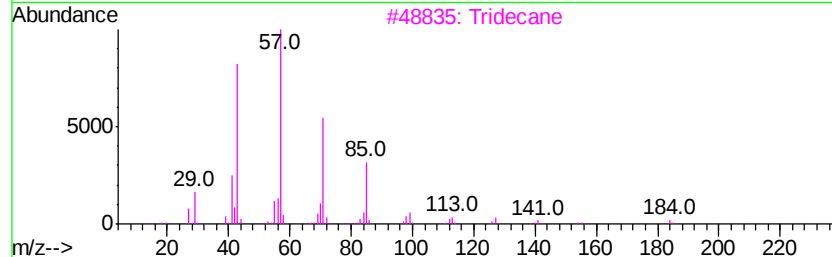
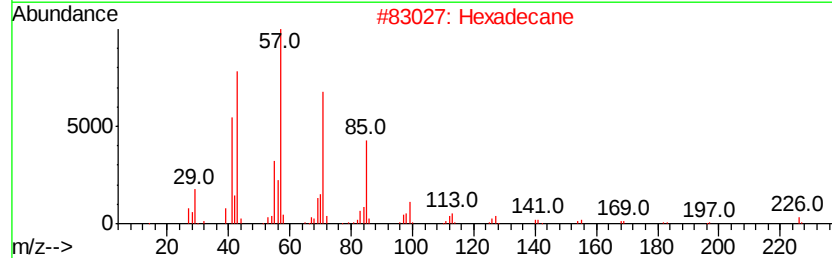
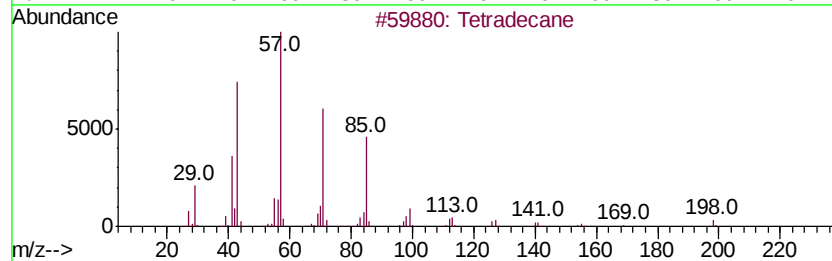
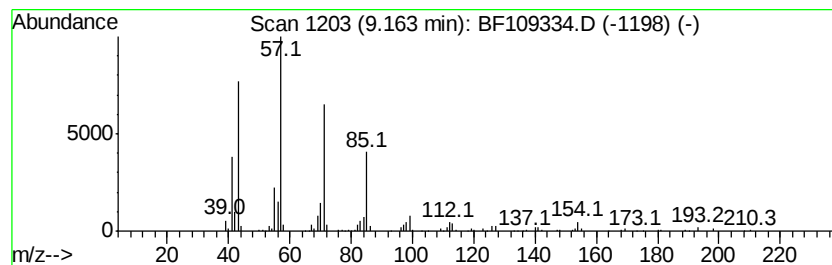
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TIC Integration Parameters: LSCINT.P

Peak Number 6 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.16	4.46 ng	545706	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	94
2		Hexadecane	226	C16H34	000544-76-3	91
3		Tridecane	184	C13H28	000629-50-5	90
4		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	78
5		Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109334.D
Acq On : 26 Sep 2018 14:17
Operator : JU/SJ
Sample : J5064-01
Misc :
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ClientSampleId :
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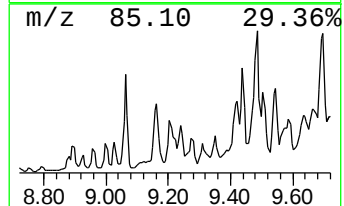
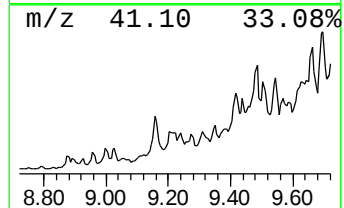
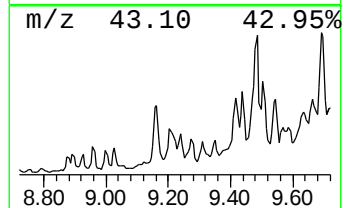
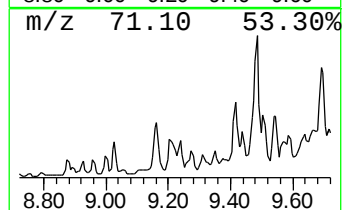
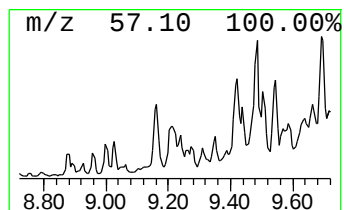
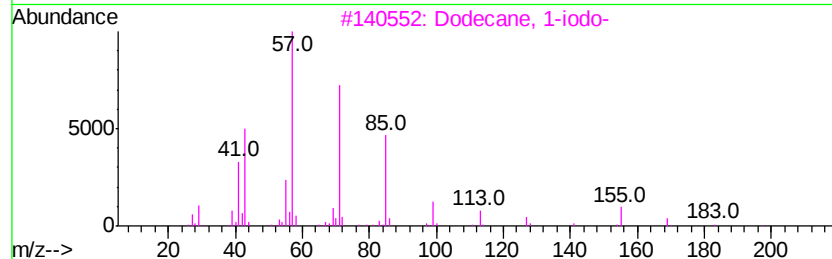
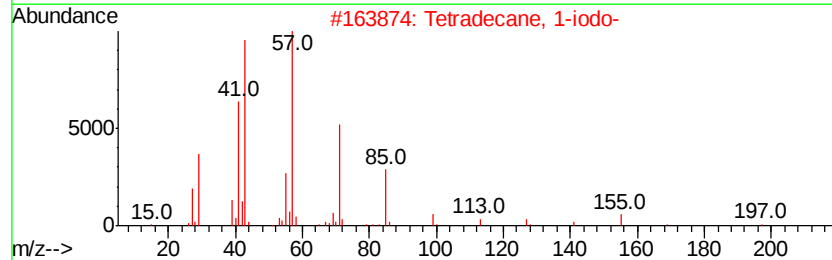
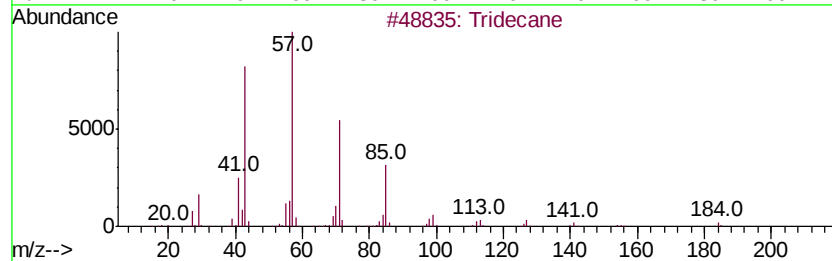
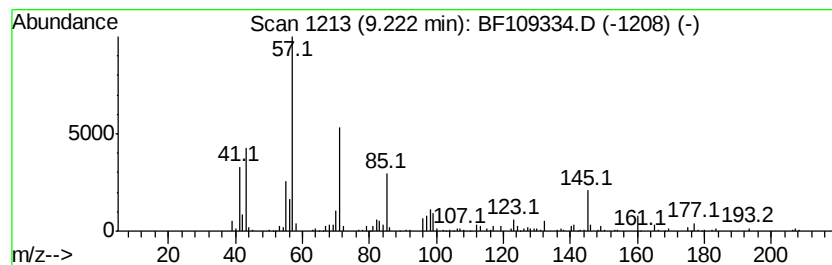
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Tridecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.22	2.74 ng	334794	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	52
2		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	50
3		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	50
4		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	50
5		Tetradecane	198	C14H30	000629-59-4	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109334.D
Acq On : 26 Sep 2018 14:17
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Sample : J5064-01
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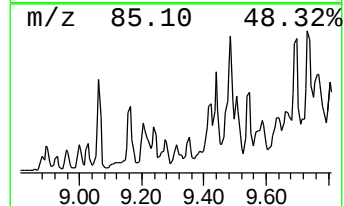
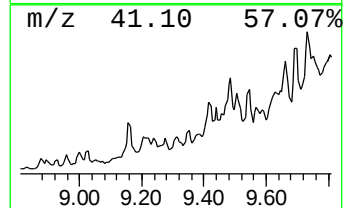
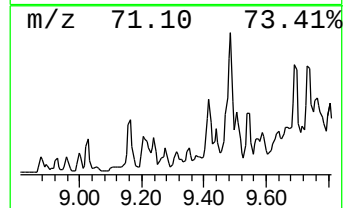
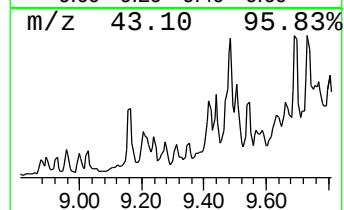
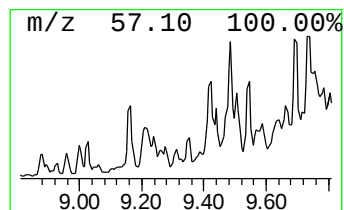
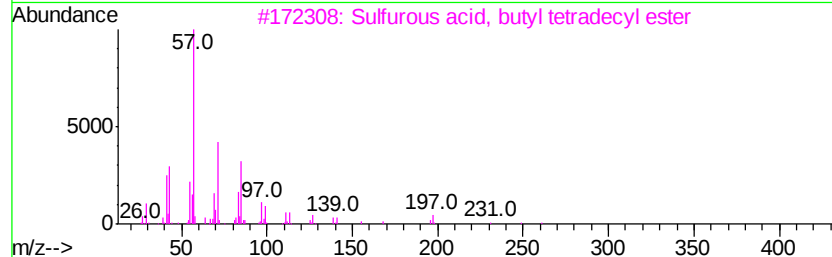
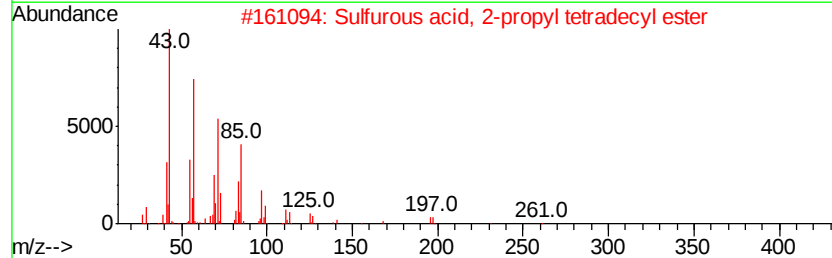
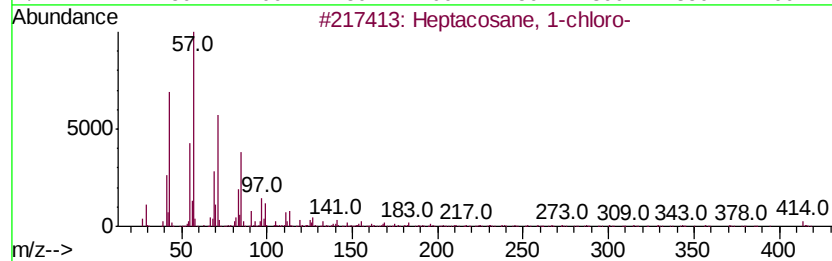
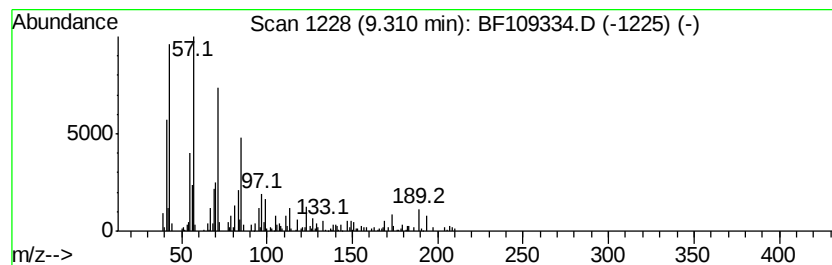
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Heptacosane, 1-chloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.31	2.48 ng	303266	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	74
2		Sulfurous acid, 2-propyl tetradecyl...	320	C17H36O3S	1000309-12-5	74
3		Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	68
4		Sulfurous acid, butyl tridecyl e...	320	C17H36O3S	1000309-18-0	68
5		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109334.D
Acq On : 26 Sep 2018 14:17
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Sample : J5064-01
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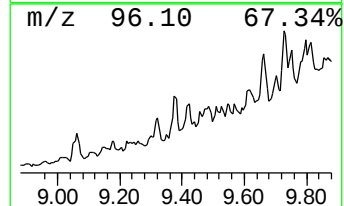
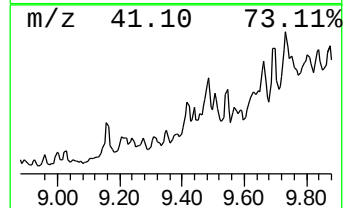
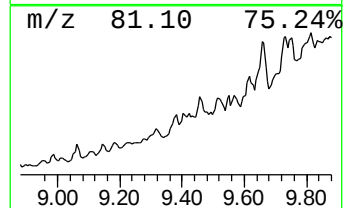
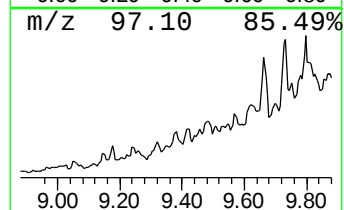
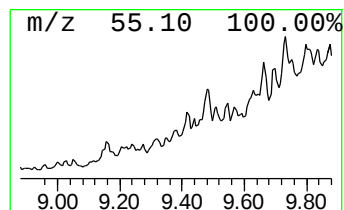
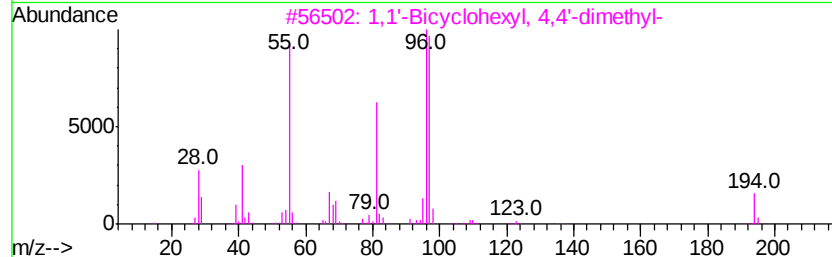
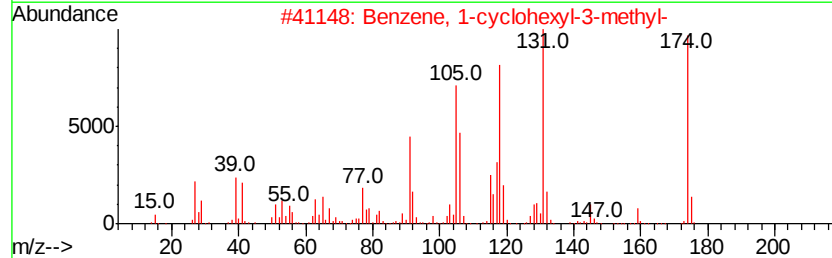
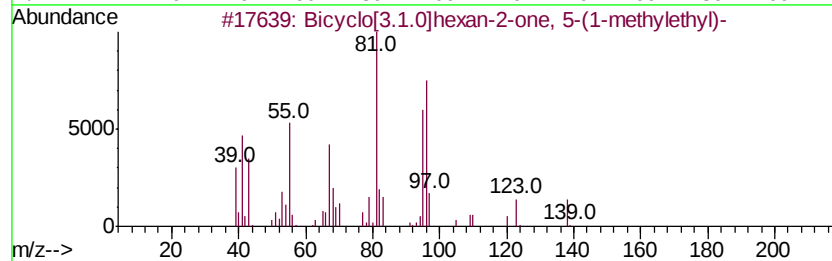
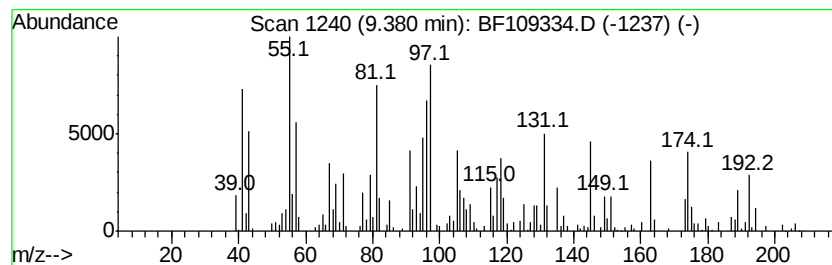
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown9.38 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.38	2.20 ng	268987	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicvclo[3.1.0]hexan-2-one, 5-(1-...	138	C9H14O	000513-20-2	38
2		Benzene, 1-cyclohexyl-3-methyl-	174	C13H18	004575-46-6	25
3		1,1'-Bicyclohexyl, 4,4'-dimethyl-	194	C14H26	054823-99-3	22
4		Cyclohexane, 1-(cyclohexylmethyl)-	194	C14H26	054823-94-8	22
5		1,3-Pentadiene, 2,3-dimethyl-	96	C7H12	001113-56-0	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109334.D
Acq On : 26 Sep 2018 14:17
Operator : JU/SJ
Sample : J5064-01
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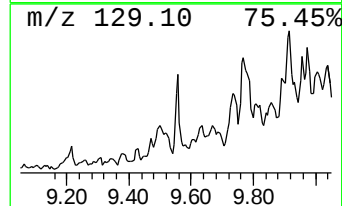
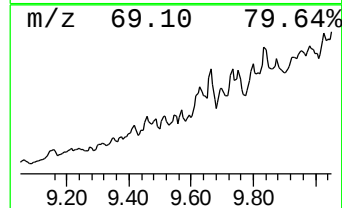
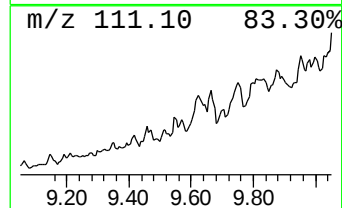
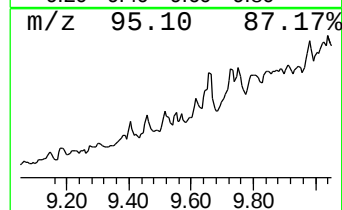
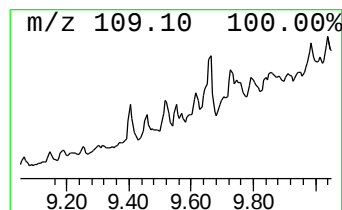
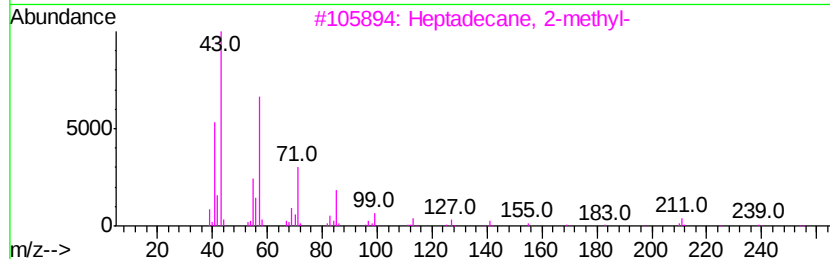
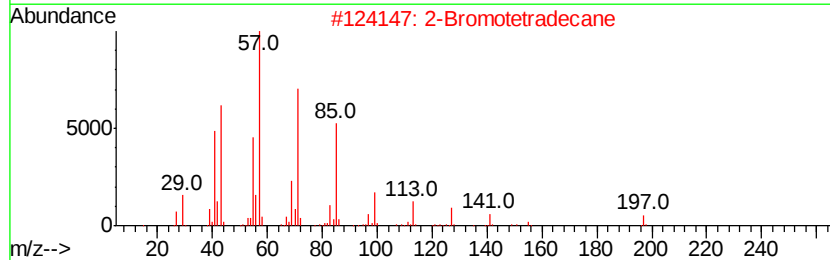
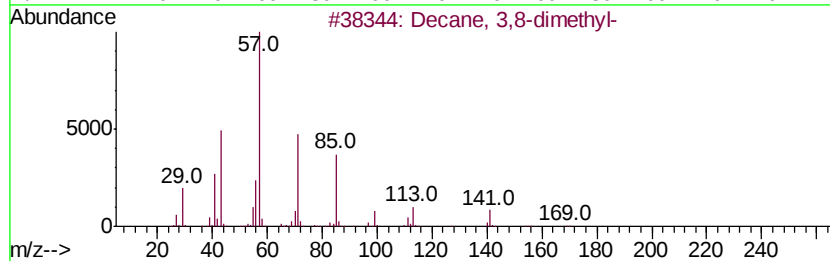
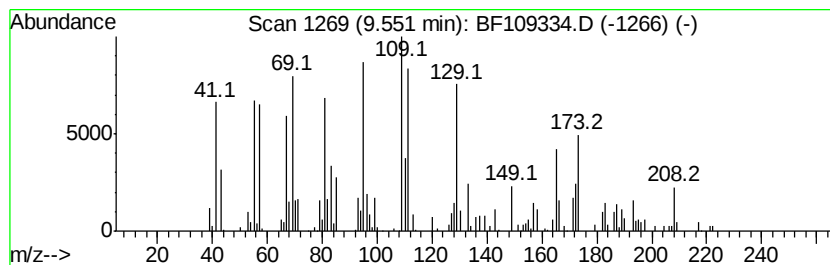
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Decane, 3,8-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.55	2.23 ng	272503	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
2		2-Bromotetradecane	276	C14H29Br	074036-95-6	58
3		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	53
4		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	53
5		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109334.D
Acq On : 26 Sep 2018 14:17
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Sample : J5064-01
Misc :
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Instrument :
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ClientSampleId :
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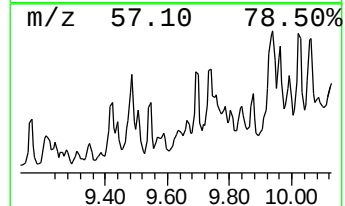
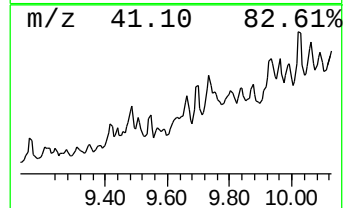
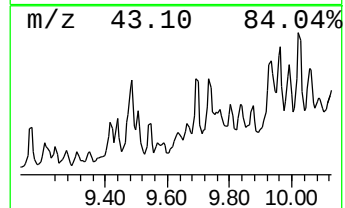
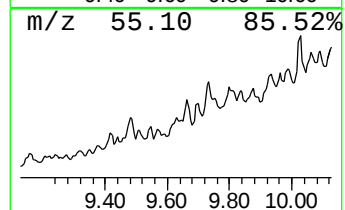
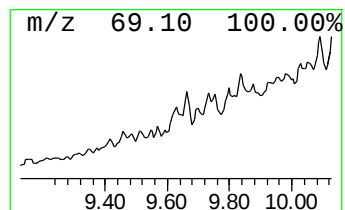
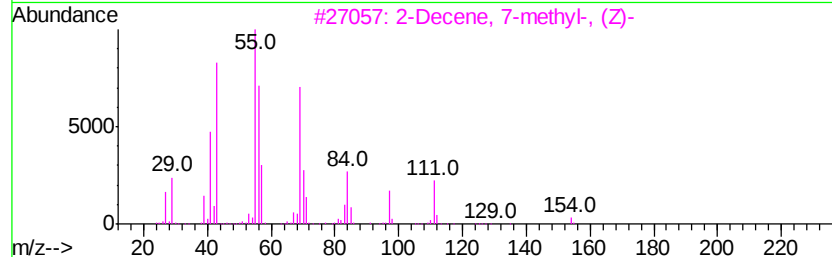
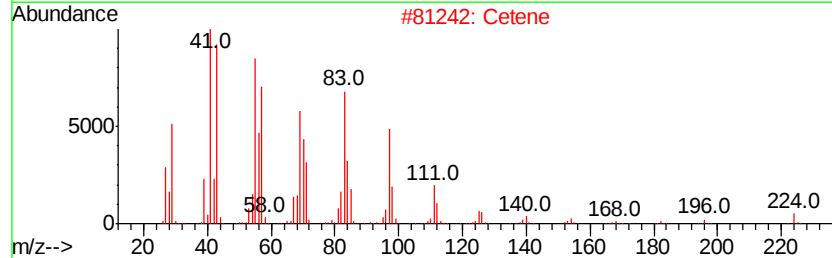
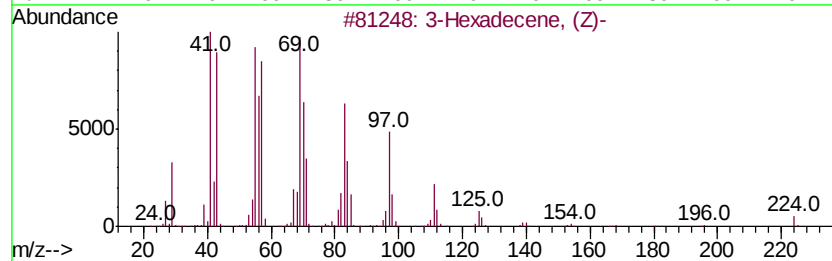
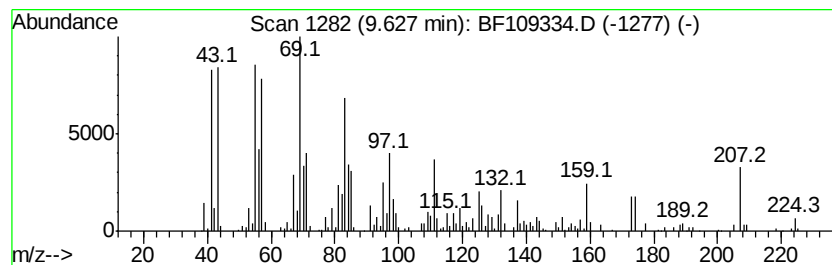
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 3-Hexadecene, (Z)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	5.55 ng	678497	Acenaphthene-d10	9.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Hexadecene, (Z)-	224	C16H32	034303-81-6	55
2			Cetene	224	C16H32	000629-73-2	47
3			2-Decene, 7-methyl-, (Z)-	154	C11H22	074630-23-2	43
4			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	43
5			Cyclohexadecane	224	C16H32	000295-65-8	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
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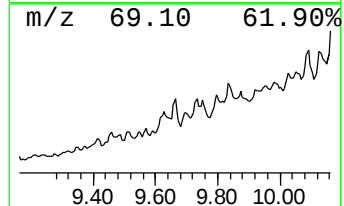
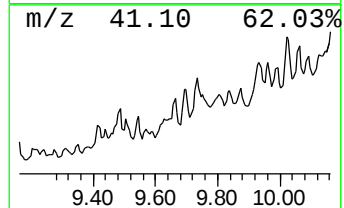
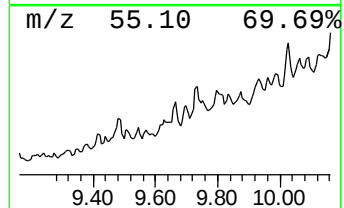
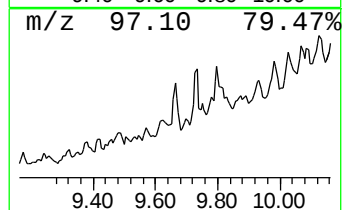
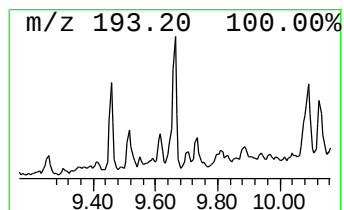
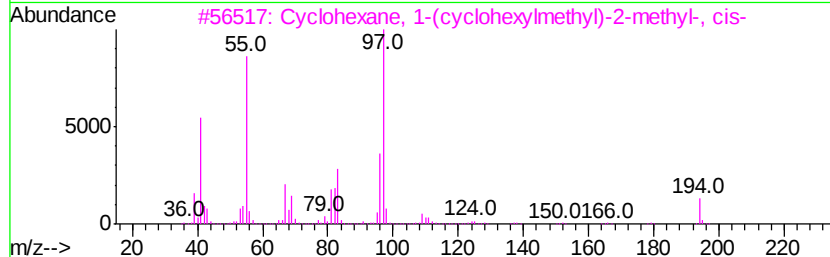
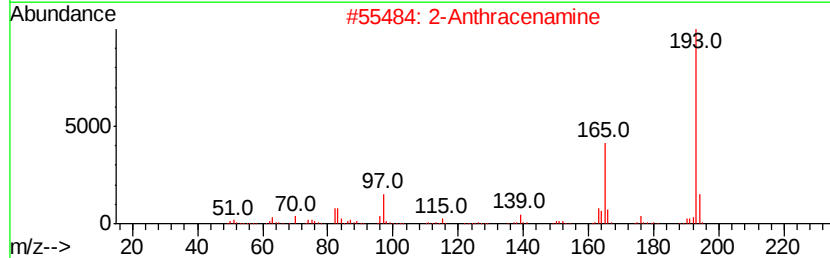
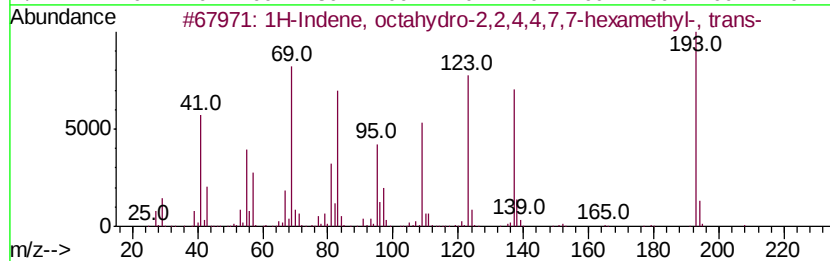
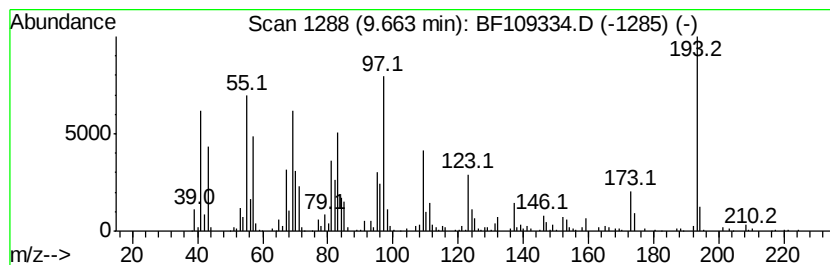
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown9.66 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.66	5.57 ng	681349	Acenaphthene-d10	9.74	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	47
2	2-Anthracenamine	193	C14H11N	000613-13-8	46
3	Cyclohexane, 1-(cyclohexylmethyl)...	194	C14H26	054824-04-3	30
4	Cyclohexane, 1-(cyclohexylmethyl)...	194	C14H26	054823-96-0	30
5	Cyclohexane, 1,5-diisopropyl-2,3...	196	C14H28	1000149-58-8	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109334.D
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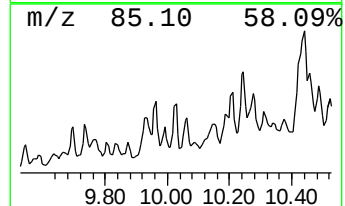
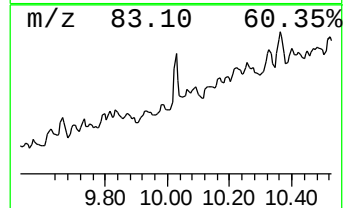
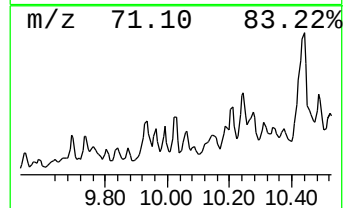
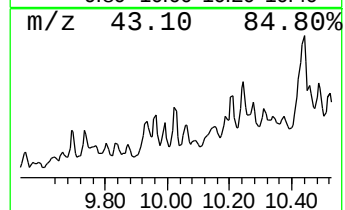
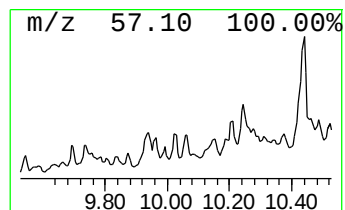
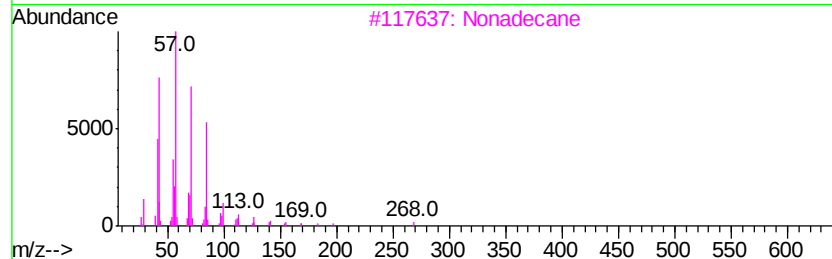
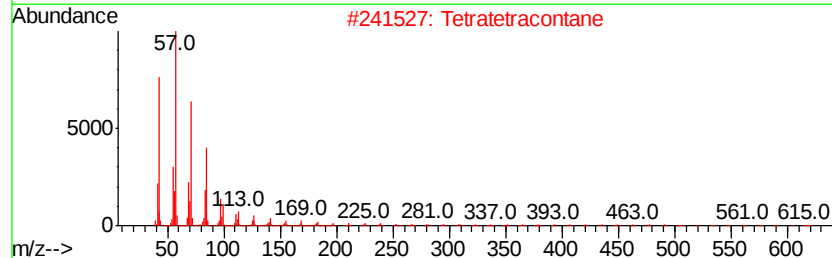
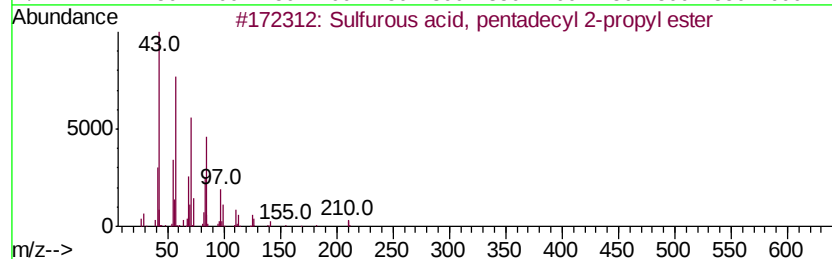
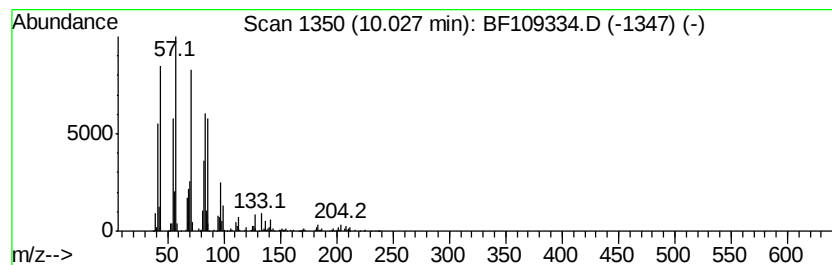
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Sulfurous acid, pentadecyl ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.03	9.70 ng	1185400	Acenaphthene-d10	9.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	64
2			Tetratetracontane	619	C44H90	007098-22-8	58
3			Nonadecane	268	C19H40	000629-92-5	58
4			Tetracosane	338	C24H50	000646-31-1	52
5			10-Methylnonadecane	282	C20H42	056862-62-5	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109334.D
Acq On : 26 Sep 2018 14:17
Operator : JU/SJ
Sample : J5064-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP

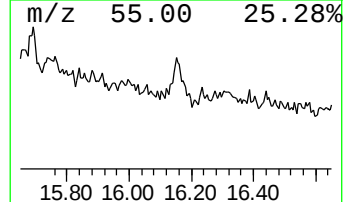
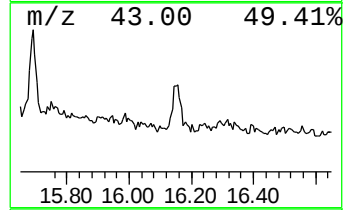
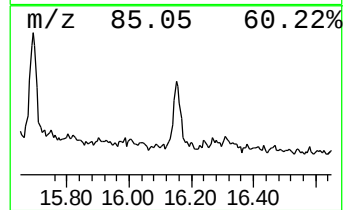
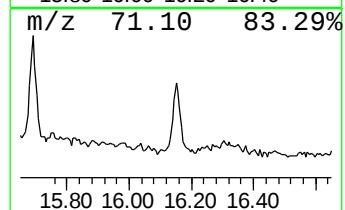
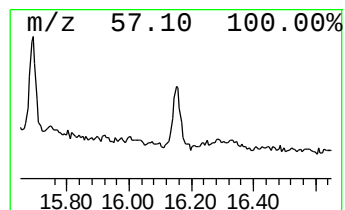
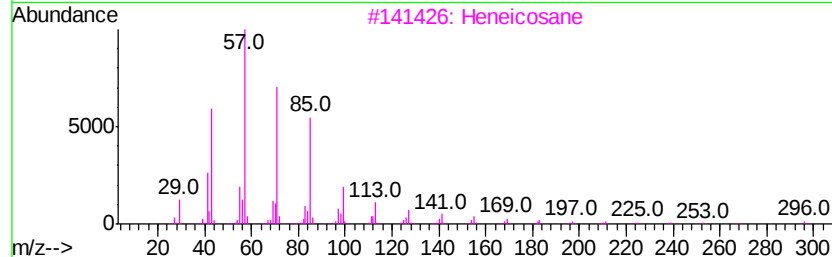
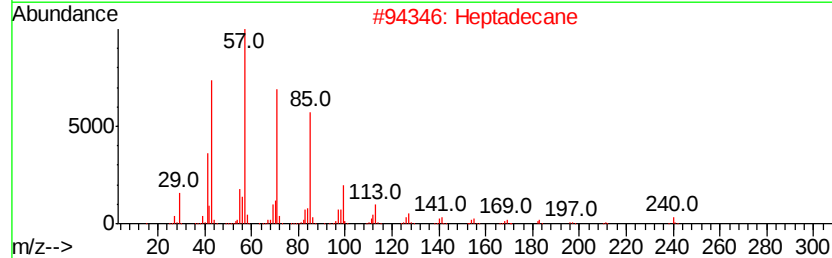
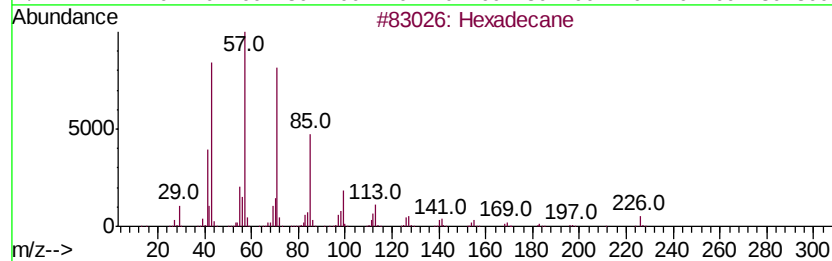
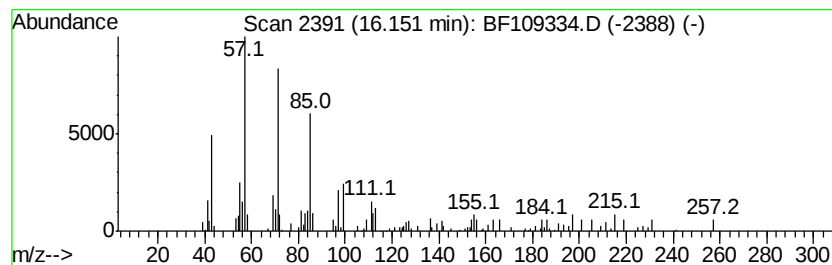
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	2.36 ng	84416	Perylene-d12	15.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	72
2		Heptadecane	240	C17H36	000629-78-7	72
3		Heneicosane	296	C21H44	000629-94-7	64
4		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	64
5		Tridecane	184	C13H28	000629-50-5	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109334.D
Acq On : 26 Sep 2018 14:17
Operator : JU/SJ
Sample : J5064-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP

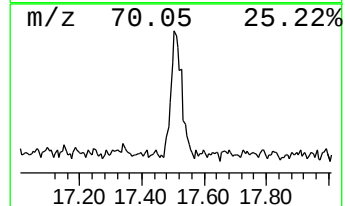
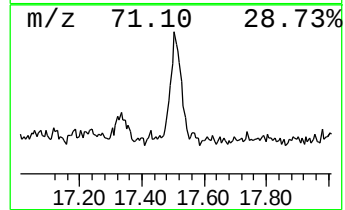
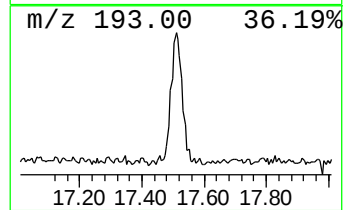
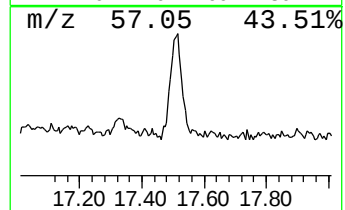
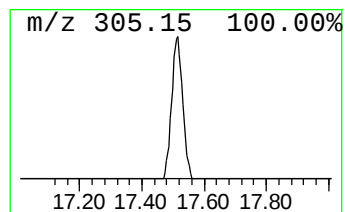
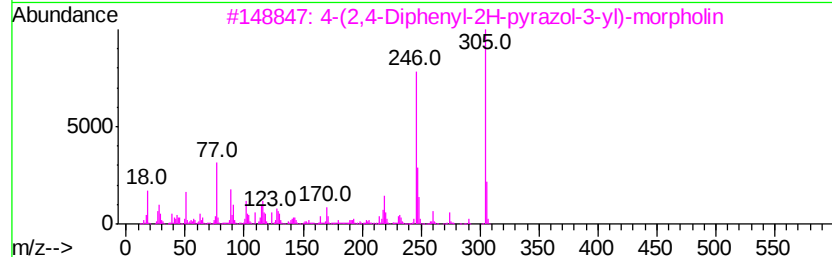
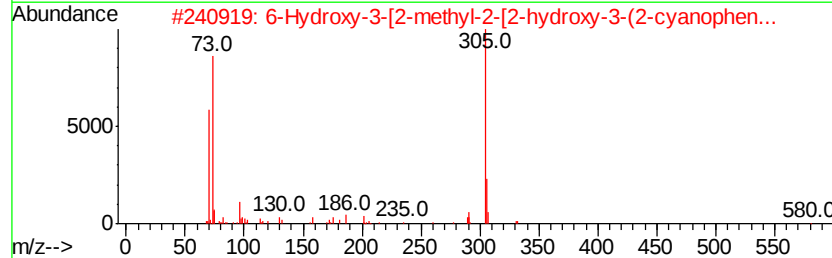
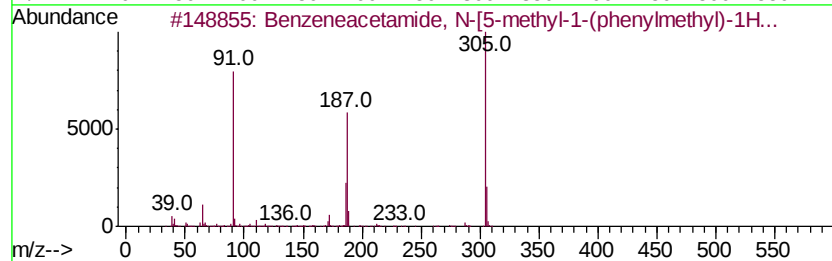
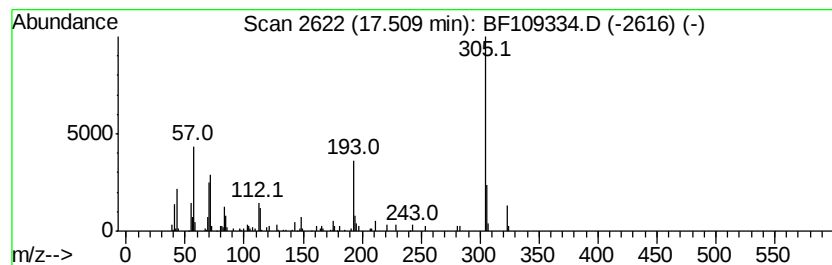
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown17.51 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.51	4.44 ng	158859	Perylene-d12	15.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneacetamide, N-[5-methyl-1-...	305	C19H19N3O	1000337-48-2	32
2		6-Hydroxy-3-[2-methyl-2-[2-hydro...	595	C31H49N3O3Si3	1000292-78-8	23
3		4-(2,4-Diphenyl-2H-pyrazol-3-yl)...	305	C19H19N3O	088743-50-4	9
4		2H-Furo[2,3-H]chromen-2-one, 8,9...	305	C19H15N3O	1000271-14-0	9
5		Furo[2,3-H]coumarine, 6-methyl-1...	305	C19H15N3O	1000270-15-3	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092618\
 Data File : BF109334.D
 Acq On : 26 Sep 2018 14:17
 Operator : JU/SJ
 Sample : J5064-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one, 4...	4.23	20.4	ng	631863	1	6.70	618632	20.0
2-Pentanone, 4-hy...	4.88	19.9	ng	614550	1	6.70	618632	20.0
unknown6.47	6.47	47.6	ng	1471440	1	6.70	618632	20.0
1-Hexanol, 2-ethyl-	6.78	22.0	ng	681220	1	6.70	618632	20.0
Tetradecane	9.16	4.5	ng	545706	3	9.74	2445250	20.0
Tridecane	9.22	2.7	ng	334794	3	9.74	2445250	20.0
Heptacosane, 1-ch...	9.31	2.5	ng	303266	3	9.74	2445250	20.0
unknown9.38	9.38	2.2	ng	268987	3	9.74	2445250	20.0
Decane, 3,8-dimet...	9.55	2.2	ng	272503	3	9.74	2445250	20.0
3-Hexadecene, (Z)-	9.63	5.5	ng	678497	3	9.74	2445250	20.0
unknown9.66	9.66	5.6	ng	681349	3	9.74	2445250	20.0
Sulfurous acid, p...	10.03	9.7	ng	1185400	3	9.74	2445250	20.0
Hexadecane	16.15	2.4	ng	84416	6	15.26	715442	20.0
unknown17.51	17.51	4.4	ng	158859	6	15.26	715442	20.0